

EAST Search History

S072 10/643,110

Ref #	Hits	Search Query	DBs	Default Operator	Plurals	Time Stamp
L1	6455	precursor\$1 same nitride	US-PGPUB; USPAT; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/06/16 15:13
L2	292	precursor\$1 same nitride same halide\$1	US-PGPUB; USPAT; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/06/16 15:13
L3	125	2 and (amine or amide)	US-PGPUB; USPAT; EPO; JPO; DERWENT; IBM_TDB	OR	ON	2006/06/16 15:13

06/16/2006 10643110 Everhart

=> d his full

(FILE 'HOME' ENTERED AT 10:21:17 ON 16 JUN 2006)

FILE 'REGISTRY' ENTERED AT 10:21:30 ON 16 JUN 2006
L1 STRUCTURE uploaded
D L1

L2 FILE 'MARPAT' ENTERED AT 10:23:09 ON 16 JUN 2006
0 SEA ABB=ON PLU=ON US2005042888/PN

L3 FILE 'REGISTRY' ENTERED AT 11:04:30 ON 16 JUN 2006
STRUCTURE UPLOADED

L4 50 SEA SSS SAM L3
L5 5629 SEA SSS FUL L3

```
FILE 'CAPLUS' ENTERED AT 11:06:28 ON 16 JUN 2006
      SET LINE 250
      SET DETAIL OFF
      E MOCVD+ALL/CT
      SET LINE LOGIN
      SET DETAIL LOGIN
```

L6 285107 SEA ABB=ON PLU=ON MOCVD OR CVD OR VAPOR DEPOS? OR ATOMIC LAYER DEPOS? OR ALD OR ALCVD OR MONOLAYER OR ATOMIC LAYER EPITAR? OR BINARY REACTION OR SEQUENTIAL SURFACE REACT? OR (PULSED OR PULSING OR PULSE) (2W) (PRECUR?)

L7 38 SEA ABB=ON PLU=ON L5 AND L6
D IBIB ABS HITSTR HITIND 1-38

FILE 'STNGUIDE' ENTERED AT 11:13:56 ON 16 JUN 2006

L8 FILE 'REGISTRY' ENTERED AT 11:19:21 ON 16 JUN 2006
L8 STRUCTURE uploaded
L8 D L8
L8 7 SEA SCS SAM LS

L10 134 SEA SSS FUL L8

FILE CATEGS ENTERED AT 11:20:28 ON 18 JUN 2008
L11 0 SEA ABB=ON PLU=ON L10 AND L6

FILED - SINGRUD - ENTERED AT 11:21:11 ON 16 JUN 2006

FILE : REGISTRY ENTERED AT 11:22:27 ON 16 JUN 2006
L12 STRUCTURE uploaded
D L12
L12 15 SEP 2006 11:22:27

L13 15 SEA SSS SAM L12
L14 341 SEA SSS FUL L12

FILE 'CAPLUS' ENTERED AT 11:23:39 ON 16 JUN 2006
L15 1 SEA ABB=ON PLU=ON L14 AND L6
L16 0 SEA ABB=ON PLU=ON L15 NOT L7.

FILE 'STNGUIDE' ENTERED AT 11:24:10 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:25:35 ON 16 JUN 2006
L17 STRUCTURE uploaded

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L18 9 SEA SSS SAM L17
L19 811 SEA SSS FUL L17

FILE 'CAPLUS' ENTERED AT 11:27:06 ON 16 JUN 2006
L20 3 SEA ABB=ON PLU=ON L19 AND L6
L21 0 SEA ABB=ON PLU=ON L20 NOT L7

FILE 'STNGUIDE' ENTERED AT 11:28:37 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:30:10 ON 16 JUN 2006
L22 STRUCTURE uploaded
L23 3086 SEA SSS FUL L22

FILE 'CAPLUS' ENTERED AT 11:30:43 ON 16 JUN 2006
L24 28 SEA ABB=ON PLU=ON L23 AND L6
L25 0 SEA ABB=ON PLU=ON L24 NOT L7

FILE 'STNGUIDE' ENTERED AT 11:32:09 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:35:33 ON 16 JUN 2006
L26 STRUCTURE uploaded
D L26
L27 0 SEA SSS SAM L26
L28 0 SEA SSS FUL L26

FILE 'STNGUIDE' ENTERED AT 11:36:32 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:38:52 ON 16 JUN 2006
L29 STRUCTURE uploaded
D L29
L30 7 SEA SSS SAM L29
L31 274 SEA SSS FUL L29

FILE 'CAPLUS' ENTERED AT 11:40:02 ON 16 JUN 2006
L32 1 SEA ABB=ON PLU=ON L31 AND L6
L33 0 SEA ABB=ON PLU=ON L32 NOT L7

FILE 'STNGUIDE' ENTERED AT 11:40:36 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:42:34 ON 16 JUN 2006
L34 STRUCTURE uploaded
L35 0 SEA SSS SAM L34
L36 0 SEA SSS FUL L34

FILE 'STNGUIDE' ENTERED AT 11:43:19 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:46:39 ON 16 JUN 2006
L37 STRUCTURE uploaded
L38 0 SEA SSS SAM L37
L39 0 SEA SSS FUL L26

FILE 'STNGUIDE' ENTERED AT 11:48:39 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 11:52:57 ON 16 JUN 2006
L40 STRUCTURE uploaded
D L40
L41 0 SEA SSS SAM L40
L42 0 SEA SSS FUL L40
L43 1 SEA ABB=ON PLU=ON 7440-33-7

06/16/2006 10643110 Everhart

D FIDE

FILE 'CAPLUS' ENTERED AT 11:57:16 ON 16 JUN 2006

FILE 'REGISTRY' ENTERED AT 13:40:35 ON 16 JUN 2006

L44 STRUCTURE uploaded
D L44

L45 50 SEA SSS SAM L44
L46 28399 SEA SSS FUL L44

FILE 'CAPLUS' ENTERED AT 13:42:39 ON 16 JUN 2006

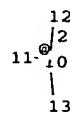
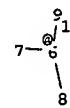
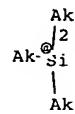
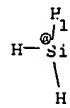
L47 2336 SEA ABB=ON PLU=ON L46 AND L6
L48 349 SEA ABB=ON PLU=ON L47 AND ?PRECURSOR?
S SILICON NITRIDE/CN

FILE 'REGISTRY' ENTERED AT 13:45:14 ON 16 JUN 2006

L49 1 SEA ABB=ON PLU=ON SILICON NITRIDE/CN

FILE 'CAPLUS' ENTERED AT 13:45:14 ON 16 JUN 2006

L50 78165 SEA ABB=ON PLU=ON L49
L51 23 SEA ABB=ON PLU=ON L48 AND L50
L52 22 SEA ABB=ON PLU=ON L51 NOT L7
D IBIB ABS HITSTR 1-22



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

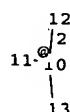
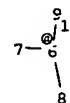
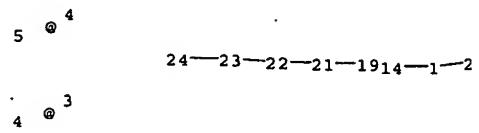
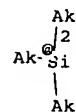
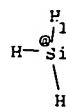
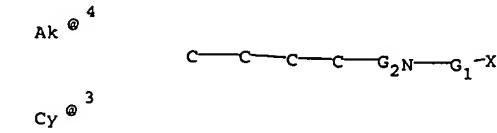
6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[*1],[*2]

G2:[*3],[*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS
13:CLASS14:CLASS19:CLASS21:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21 22 23 24

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21 21-22 22-23 23-24

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

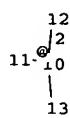
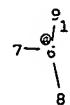
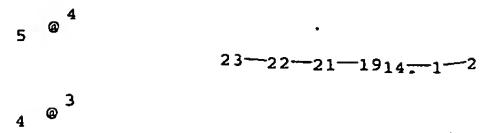
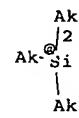
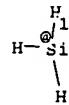
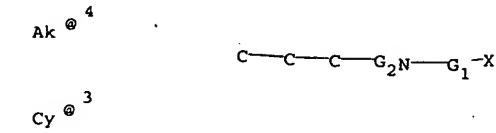
6-7 6-8 6-9 21-22 22-23 23-24

G1:Si,B,Al,Nb,Ta,Ti,W,[*1],[*2]

G2:[*3],[*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS
 13:CLASS14:CLASS19:CLASS21:CLASS22:CLASS23:CLASS24:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21 22 23

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21 21-22 22-23

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

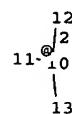
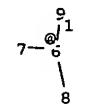
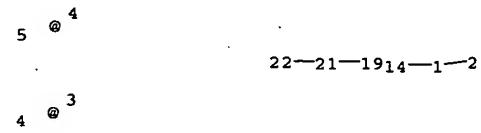
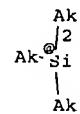
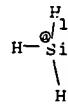
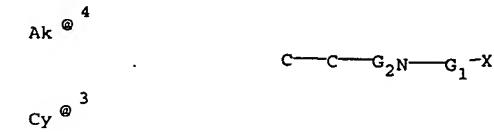
6-7 6-8 6-9 21-22 22-23

G1:Si,B,Al,Nb,Ta,Ti,W,[*1],[*2]

G2:[*3],[*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS
 13:CLASS14:CLASS19:CLASS21:CLASS22:CLASS23:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21 22

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21 21-22

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

6-7 6-8 6-9 21-22

G1:Si,B,Al,Nb,Ta,Ti,W,[*1],[*2]

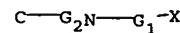
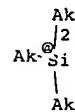
G2:[*3],[*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS14:CLASS19:CLASS21:CLASS22:CLASS

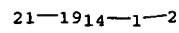
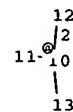
Ak @ 4

Cy @ 3



5 @ 4

4 @ 3



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19 21

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21

exact bonds :

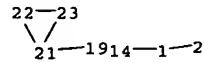
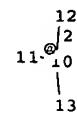
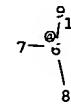
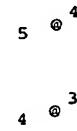
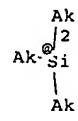
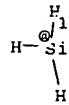
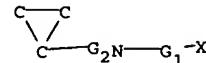
6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[*1],[*2]

G2:[*3],[*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS
13:CLASS14:CLASS19:CLASS21:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19

ring nodes :

21 22 23

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

ring bonds :

21-22 21-23 22-23

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21 21-22 21-23 22-23

exact bonds :

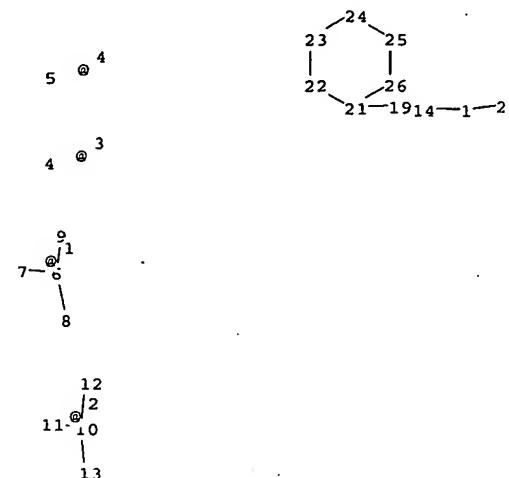
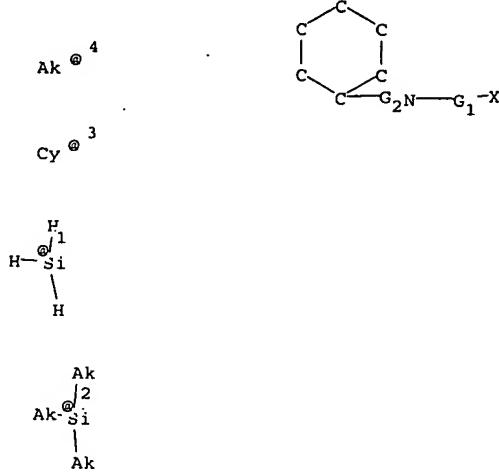
6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[*1],[*2]

G2:[*3],[*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS
13:CLASS14:CLASS19:CLASS21:CLASS22:CLASS23:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19

ring nodes :

21 22 23 24 25 26

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

ring bonds :

21-22 21-26 22-23 23-24 24-25 25-26

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21 21-22 21-26 22-23 23-24 24-25 25-26

exact bonds :

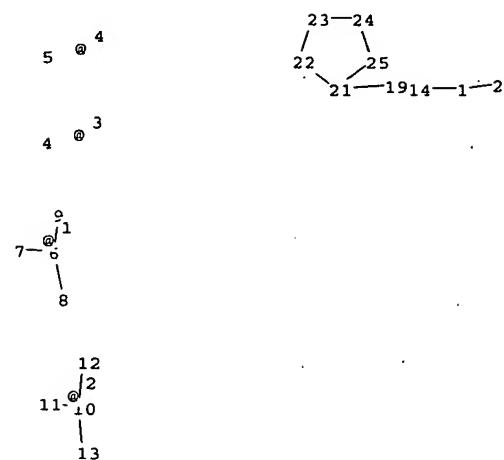
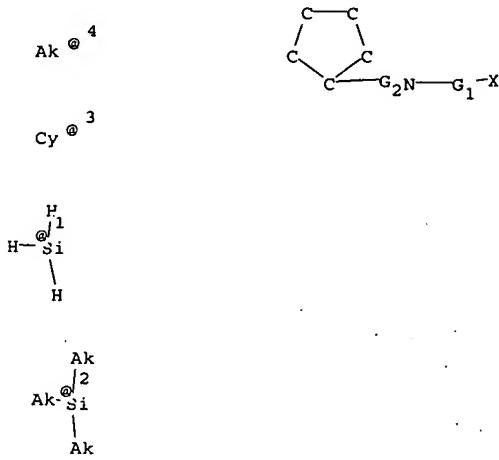
6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[*1],[*2]

G2:[*3],[*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS14:CLASS19:CLASS21:Atom 22:CLASS23:CLASS24:CLASS25:Atom 26:Atom



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19

ring nodes :

21 22 23 24 25

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

ring bonds :

21-22 21-25 22-23 23-24 24-25

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21 21-22 21-25 22-23 23-24 24-25

exact bonds :

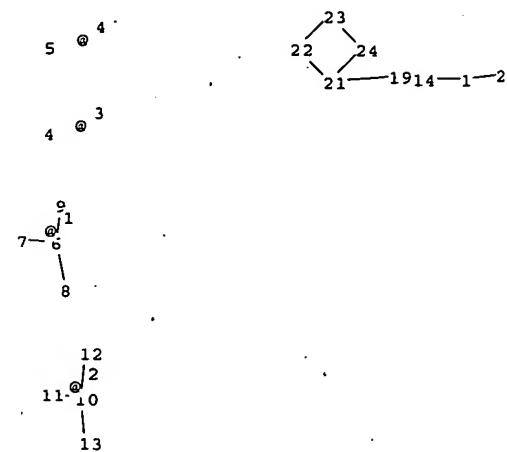
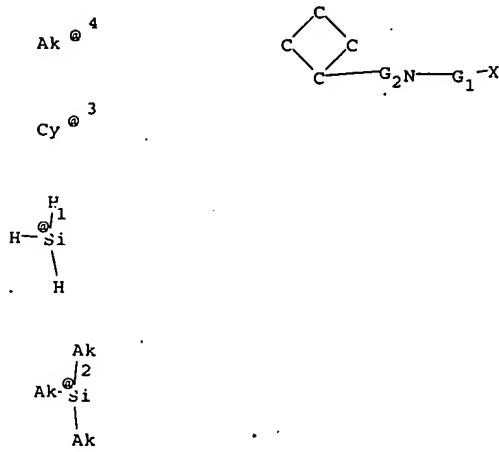
6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[*1],[*2]

G2:[*3],[*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS14:CLASS19:CLASS21:Atom 22:Atom 23:Atom 24:Atom 25:CLASS



chain nodes :

1 2 4 5 6 7 8 9 10 11 12 13 14 19

ring nodes :

21 22 23 24

chain bonds :

1-2 1-14 6-7 6-8 6-9 10-11 10-12 10-13 14-19 19-21

ring bonds :

21-22 21-24 22-23 23-24

exact/norm bonds :

1-2 1-14 10-11 10-12 10-13 14-19 19-21 21-22 21-24 22-23 23-24

exact bonds :

6-7 6-8 6-9

G1:Si,B,Al,Nb,Ta,Ti,W,[*1],[*2]

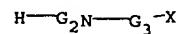
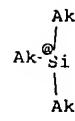
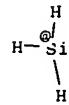
G2:[*3],[*4]

Match level :

1:CLASS2:CLASS4:Atom 5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS14:CLASS19:CLASS21:Atom 22:Atom 23:Atom 24:CLASS

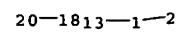
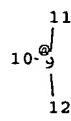
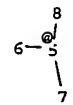
Ak @ 2

Cy @ 1



4 @ 2

3 @ 1



chain nodes :

1 2 3 4 5 6 7 8 9 10 11 12 13 18 20

chain bonds :

1-2 1-13 5-6 5-7 5-8 9-10 9-11 9-12 13-18 18-20

exact/norm bonds :

1-2 1-13 9-10 9-11 9-12 13-18 18-20

exact bonds :

5-6 5-7 5-8

G2:[*1],[*2]

G3:Ta,Ti

Match level :

1:CLASS2:CLASS3:Atom 4:CLASS5:CLASS6:CLASS7:CLASS8:CLASS9:CLASS10:CLASS11:CLASS12:CLASS13:CLASS18:CLASS20:CLASS



chain nodes :

1 3 4 7 9 10

chain bonds :

1-10 7-9 7-10

exact/norm bonds :

7-9 7-10

exact bonds :

1-10

G1:Si,B,Al,Nb,Ta,Ti,W

G2:[*1],[*2]

Match level :

1:CLASS3:Atom 4:CLASS7:CLASS9:CLASS10:CLASS